ANALYTICAL RESULTS OF SURFACE WATER SAMPLES COLLECTED FROM RACCOON CREEK

July 22, 1999 Sampling Event

Prepared for

Lyondell Chemical Worldwide Inc.Beazer East, Inc.

Prepared by

Applied Hydrology Associates, Inc.
Pittsburgh, PA
Denver, CO

August 18, 1999



TABLE OF CONTENTS

Section	<u>P:</u>	age
1.0	INTRODUCTION	1
2.0	SAMPLING	1
3.0	RESULTS	2
Tables		
1 2	Summary of Analytical Results for Samples Collected from Raccoon Creek Historic Benzene Concentrations at Transect E	2 3
<u>Figure</u>	<u>s</u>	
1 2	Transect Location Map Surface Water Benzene Concentrations at Transect E	4 5
Appen	<u>dices</u>	
A R	Groundwater Elevations, East and West Sides of Raccoon Creek Data Validation Report	

1.0 INTRODUCTION

This report presents the results of surface water samples collected from Raccoon Creek at the Lyondell Chemical Worldwide, Inc. / Beazer East Inc. Monaca, PA site during the July 22, 1999 quarterly monitoring event. The samples were collected in compliance with Appendix D of the 1997 Consent Order and Agreement (1997 CO&A) between ARCO Chemical Company¹, BEI and the Pennsylvania Department of Environmental Protection (PADEP) dated October 20, 1997.

2.0 SAMPLING

Surface water samples were collected at Transect E as defined in the 1997 CO&A. The location of Transect E is shown in Figure 1. In addition, water elevations were measured in nearby monitoring wells and the results are presented in Appendix A.

A total of eight surface water samples, including a duplicate were collected from Raccoon Creek on July 22, 1999. These samples were collected at the same three locations along Transect E as in previous sampling events. The locations are shown in Figure 2 and are at the center of the stream, and approximately 30 feet from the east and west banks. At the center location, samples were collected at three depths; 6 inches below the surface, 2 inches above the bottom, and midway between the surface and bottom. Samples from the east and west sides of the transect were collected at two depths; 2 inches above the bottom, and midway between the surface and bottom.

During sampling a boat was stationed at Transect E using a rope secured to the east and west shores of Raccoon Creek. The samples were collected by using a peristaltic pump to pump water from the desired depth into three 40-ml vials preserved with hydrochloric acid. Samples were collected from the required depths utilizing tubing secured to a vertical steel rod lowered from the boat until it rested on the bottom of the creek. The rod did not penetrate the sediments on the creek bottom because a 1-foot diameter disc constructed of steel mesh is fastened perpendicular to the bottom of the rod.

Two tubes were used. The bottom of the "deep sample tube" was secured to the probe 2 inches from the bottom of the probe. The bottom of the "mid-depth sample tube" is adjustable and was secured to the probe mid-depth at each location. Care was taken not to disturb the sediments at the sampling location and the pumped water was closely monitored to ensure sediment was not included in the sample. One gallon of water was pumped through the tubing before each sample was obtained in order to purge the tubing.

¹ ARCO Chemical Company is now Lyondell Chemical Worldwide

The samples were uniquely numbered as follows to identify the location, depth and date of sampling:

RC-EC-00-0799

Where:

RC indicates Raccoon Creek

EC indicates Transect E and location (C = center, L = left bank, R = right bank [facing downstream])

00 indicates sample depth in feet and tenths of a foot (0.0 feet)

0799 indicates the month and year collected (July 1999)

Samples were logged onto a chain of custody form (included in of the Analytical Report in Appendix B) and stored on ice until receipt by Reliance Laboratories Inc. in Edison, NJ. Reliance analyzed the samples using USEPA Method 524.2 for BTEXS.

3.0 RESULTS

The analytical results are presented in Table 1. No BTEXS constituents were detected in any of the eight samples. Sampling locations and depths are shown on Figure 2, along with the concentration of benzene at each location. Water levels in wells near Raccoon Creek are presented in Appendix A.

Table 1
Summary of Analytical Results for Samples Collected from Raccoon Creek

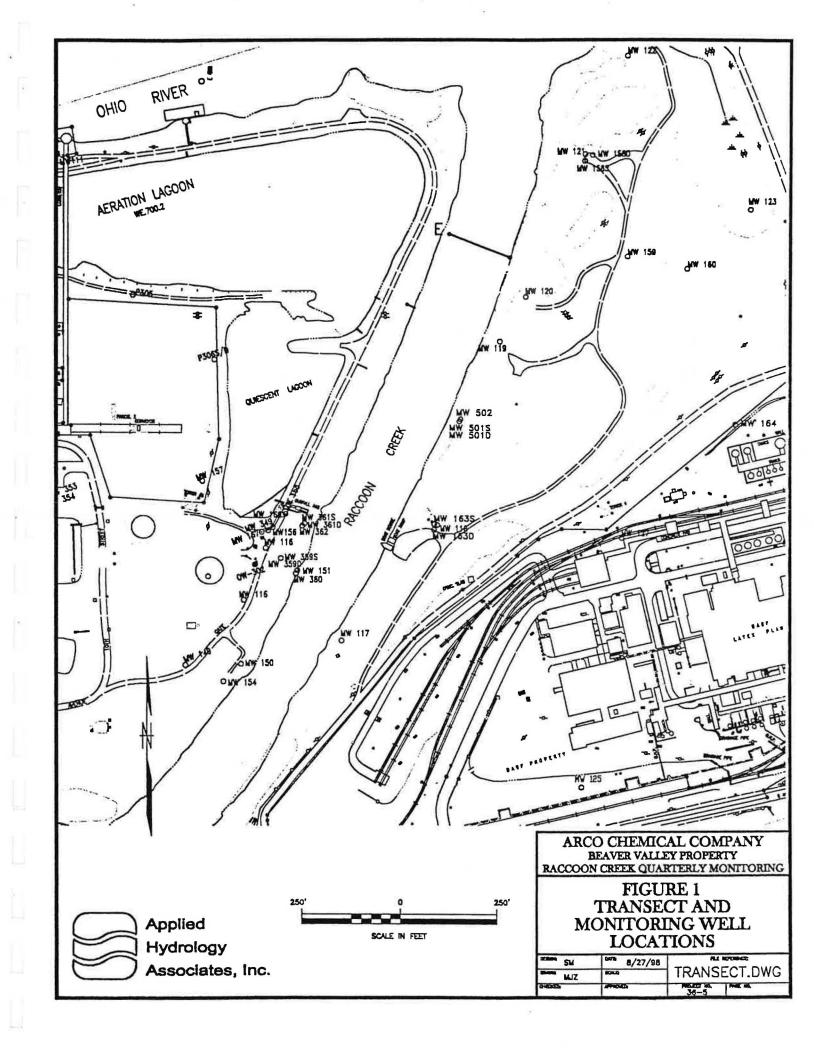
Sample Name	Benzene	Toluene	Ethylbenzene	Xylene	Styrene
RC-EL-24-0799	< 0.13	< 0.6	< 0.22	< 0.22	< 0.58
RC-EL-12-0799	< 0.13	< 0.6	< 0.22	< 0.22	< 0.58
RC-EC-00-0799	< 0.13	< 0.6	< 0.22	< 0.22	< 0.58
RC-EC-00-0799A	< 0.13	< 0.6	< 0.22	< 0.22	< 0.58
RC-EC-35-0799	< 0.13	< 0.6	< 0.22	< 0.22	< 0.58
RC-EC-71-0799	< 0.13	< 0.6	< 0.22	< 0.22	< 0.58
RC-ER-34-0799	< 0.13	< 0.6	< 0.22	< 0.22	< 0.58
RC-ER-68-0799	< 0.13	< 0.6	< 0.22	< 0.22	< 0.58

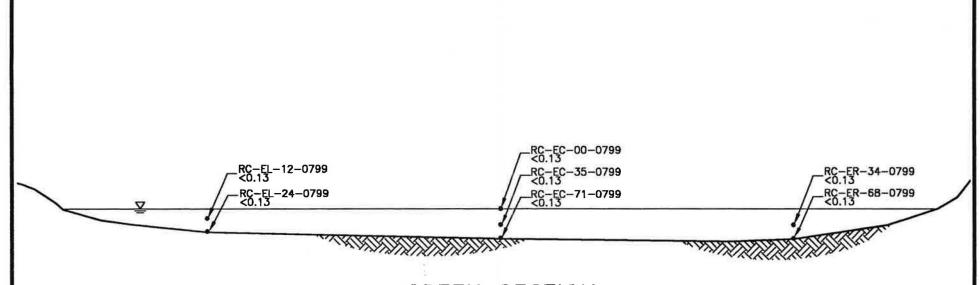
The analytical data were validated upon receipt and found to be acceptable. A Data Validation Report is provided in Appendix B. Table 2 presents the historical concentration of benzene in Raccoon Creek at Transect E during all monitoring events to date.

Historic Benzene Concentrations at Transect E (ug/L)

Sampling Location	Sampling Depth	7/23/97	10/28/97	2/25/98	5/21/98	7/29/98	10/27/98	2/3/99	4/27/99	7/22/99
30 Feet off	Mid-depth	0.28	<0.13	<0.13	0.70	<0.13	1.57(1)	0.37	< 0.66	<0.13
West Bank										
30 Feet off	Deep	0.81	<0.13	<0.13	0.70	<0.13	0.61(1)	0.49	< 0.66	<0.13
West Bank						1				
Center of	Shallow	0.24	<0.13	0.38	0.70	<0.13	<0.13	0.61 ⁽¹⁾	< 0.66(1)	<0.13(1)
Creek										
Center of	Mid-	0.18	<0.13	0.49	0.64	<0.13	0.2	0.64	< 0.66	<0.13
Creek	Depth									
Center of	Deep	0.46	<0.13	0.30	0.60	<0.13	<0.13	0.69	< 0.66	<0.13
Creek										
30 Feet off	Mid-depth	0.16	<0.13	<0.13	<0.13	0.13	0.52	< 0.13	< 0.66	<0.13
East Bank							1			
30 Feet off	Deep	<0.13	<0.13	0.14	0.22	0.22	<0.13	< 0.13	< 0.66	<0.13
East Bank										

(1) Results shown are the average of the blind duplicate samples.





CREEK SECTION LOOKING DOWNSTREAM

LEGEND

SURFACE WATER SAMPLE LOCATION
 ALL CONCENTRATIONS IN ug/L



Applied Hydrology Associates, Inc.



LYONDELL CHEMICAL WORLDWIDE BEAVER VALLEY PROPERTY RACCOON CREEK QUARTERLY MONITORING

FIGURE 2

SURFACE WATER
BENZENE CONCENTRATIONS
AT TRANSECT 'E'

JULY 22, 1999

SM	PATE 8/17/98	FLE REFERENCE
JLS	NOT TO SCALE	BENZENE.dwg
Dec. all	ATTIONES.	700.000 MA. PARK NO. 38-5

Appendix A

Groundwater Elevations, East and West Sides of Raccoon Creek

GROUNDWATER LEVELS ON THE EAST AND WEST SIDES OF RACCOON CREEK July 22, 1999

Well Number	Top of Casing (TOC) (ft. amsl)	Depth to SPL from TOC (2) (ft. amsl)	Depth to Water from TOC (2) (ft. amsl)	Calculated Water Level Elevation (1) (ft. amsl)	Calculated SPL Thickness (3) (ft. amsl)	Comments
		N	Ionitoring V	Vells Screened	in Silty Clay	Unit
MW - 360	COE 04	ND	2.39	OTH AREA 683.45	N/A	
MW - 300 MW - 170	685.84 706.70	ND ND	22.44	684.26	N/A N/A	
MW - 1/0	689,43	ND	5.91	683.52	N/A	
WI W - 302	069,43	ND		RACCOON CREEK		
			·	MCCOON CREEK	AREA	Monitoring well is dry. Screened above water table
MW- 118	690,39	ND	7.00	683,39	N/A	Bottem of well is 7.00 feet below TOC.
MW - 502	701.86	ND	18.62	683.24	N/A	Bottem of wen is 7:00 feet below 100.
MW - 119	705.59	ND	22.32	683.27	N/A	
MW - 120	709.42	ND	26.10	683.32	N/A	
MW - 121	713.90	ND	30.58	683.32	N/A	
MW - 152	696.35	ND	13.06	683.29	N/A	
14144 102	070.55					C111-'4
		Monitor	ing wens so	creened in Upp OTH AREA	er Sand and	Gravei Unit
MW - 344	709.42	ND	25.63	683.79	N/A	
MW - 359S	692,93	ND	9.46	683.47	N/A	
MW - 361S	689.40	ND	6.00	683.40	N/A	
14144 - 2012	002.40	ND	0.00	083,40	IVA	Well had excessive pressure build up from OTH
MW - 169	707.93	ND	28,58	679.35	N/A	Sparge event (7/13/99 - 7/14/99).
MW - 167	711.06	ND	27.57	683.49	N/A	Top of casing changed from 707.36 to 711.06 on 11/98 accommodate respiration monitoring well head. Monitoring well stick up is 3.70 above orig. TOC
				RACCOON CREEK		
MW - 163S	690.87	ND	7.56	683.31	N/A	
MW - 501S	701.30	ND	18.31	682.99	N/A	
MW - 162S	706.05	ND	22.80	683.25	N/A	
MW - 159	708.99	ND	25.71	683.28	N/A	
MW - 160	701.00	ND	17.69	683.31	N/A	
MW - 158S	713.60	ND	30.34	683.26	N/A	
MW - 122	692.78	ND	9.55	683.23	N/A	
Note: See figure						
				to Water from TOC.		
				e. ND means no SPL		s not applicable, no SPL was detected.

GROUNDWATER LEVELS ON THE EAST AND WEST SIDES OF RACCOON CREEK July 22, 1999

Well Number	Top of Casing (TOC) (ft. amsl)	Depth to SPL from TOC (2) (ft. amsl)	(-	Calculated Water Level Elevation (1) (ft. amsl)	Calculated SPL Thickness (3) (ft. amsl)	Comments		
		Monitori	ng Wells Sc	reened in Lowe	er Sand and	Gravel Unit		
MW 345	708.91	ND	25.60	683,31	N/A			
MW 361D	689.35	ND	5.91	683,44	N/A			
MW 359D	692.80	ND	9.43	683,37	N/A			
WW JSJD	072.00	ND		ACCOON CREEK				
MW 163D	689.62	ND	6.25	683,37	N/A			
MW 501D	701.44	ND	18.17	683.27	N/A			
MW 166D	703.95	ND	20.64	683.31	N/A			
MW 158D	712.04	ND	28.86	683.18	N/A			
		W	ater Levels i	n Raccoon Cre	ek and Ohio	River		
			RACCOO	N CREEK AREA S	TAFF GAUGE			
Time of	Staff Gauge Elevation (ft. amsl)	Staff Gauge	Calculated Water Level Elevation		Comments			
Observation	(4) (5)	Reading	(ft. amsl)					
10:25	685.00	1.80	683.80					
11:24	685.00	1.80	683.80					
			ОН	IO RIVER. STAFF	GAUGE			
7:56	685.96	3.10	683.06					
11:13	685.96	3.12	683.08			1)		
						K		
Note: See figure	1							
		Elevation of To	C minus Denth	to Water from TOC.		1		
				e. ND means no SPL	was detected			
		0				not applicable, no SPL was detected.		
			on staff gauge at					
			on staff gauge at					

Appendix B Data Validation Report



1200 South Parker Road, Suite 100

Denver, CO 80231

Tel: (303) 873-0164

Fax: (303) 873-6110

MEMORANDUM

TO:

Files

FROM:

Skip Meier, Applied Hydrology Associates

DATE:

August 17, 1999

SUBJECT:

Data Validation Results, Lyondell Chemical Worldwide Beaver Valley Property

Data validation was performed on the volatile organic analytical data from eight surface water samples obtained from Raccoon Creek on July 22, 1999 and also on a Rinsate Blank and Trip Blank. The validation was performed in accordance with the "USEPA Contract Laboratory Program National Functional Guidelines for Organic Data Review". Reliance Laboratories Inc. performed the analysis using EPA Method 524.2. The samples reviewed included:

Field Sample ID	Lab Sample ID
RC-EL-24-0799	R-6267.3
RC-EL-12-0799	R-6267.4
RC-EC-00-0799	R-6267.7
RC-EC-00-0799A	R-6267.8
RC-EC-35-0799	R-6267.6
RC-EC-71-0799	R-6267.5
RC-ER-34-0799	R-6267.2
RC-ER-68-0799	R-6267.1
Rinsate Blank	R-6267.9
Trip Bland	R-6267.10

Items reviewed and actions taken were as follows:

√ Method:

The ten samples were analyzed for BTEXS by method USEPA 524.2 on July 23, 1999.

√ Holding Time:

All Samples were analyzed within the 14-day holding time.

√ Blanks:

No target compounds were detected in the associated method blank.

√ System Monitoring Compounds:

The "USEPA Contract Laboratory Program National Functional Guidelines for Organic Data Review" indicate that "Recoveries for system monitoring compounds in volatile samples and blanks must be within the limits specified in the Method." However, Method 524.2 does not specify a required

recovery. Nevertheless, 4-bromofluorobenzene and 1,2-dichlorobenzene-d4 surrogate recoveries were within 96-122 percent and this is acceptable.

√ Internal Standards:

All fluorobenzene internal standards were within the established criteria for area internal standard and retention time.

√ GC/MS Instrument Performance Check:

All bromofluorobenzene (BFB) tunes met the ion abundance criteria. Analysis of the instrument performance check solution was performed at the beginning of each 12-hr period during which the samples were analyzed.

√ Initial Calibrations:

The initial calibration performed on July 23, 1999 for Instrument HP5971A met the 30 percent relative standard deviation (RSD) and 0.05 minimum relative response factor criteria for all compounds.

√ Continuing Calibrations:

Continuing calibration was run and compared to the correct initial calibration. All continuing calibrations met the 25 percent difference and minimum relative response factor criteria for all compounds.

√ Matrix Spike/Duplicate:

The matrix spike/duplicate results for recovery and RPD were within the Quality Control limits.

√ Target Compound Indentification/Quantitation:

No problems were identified with compound identification or quantities.

√ Field Duplicate:

A field duplicate was collected during this sampling event. The duplicate sample was denoted by an "A" at the end of the sample name. The pair is RC-EC-00-0799 and duplicate RC-EC-00-0799A. Table 1 below summarizes the RPD for the sample/duplicate pair, indicating that no BTEXS compounds were detected for either the primary sample or the duplicate.

Table 1: Relative Percent Difference (RPD)

Sample Name	Benzene (ppb)	RPD (%)	Toluene (ppb)	RPD (%)	Ethyl- Benzene (ppb)	RPD (%)	Xylene (ppb)	RPD (%)	Styrene (ppb)	RPD (%)
RC-EC-00-0799	ND	NA	ND	NA	ND	NA	ND	NA	ND	NA
RC-EC-00-0799A	ND	NA	ND	NA	ND	NA	ND	NA	ND	NA

ND = Non Detect

NA = Not Applicable

√ Summary:

No inconsistencies were noted. No BTEXS constituents were detected in the duplicate sample pair RC-EC-00-0799 and RC-EC-00-0799A (See Table 1). No BTEXS compounds were detected in either the trip blank or the field blank.

RELIANCE LABORATORIES, INC.



175 MAY STREET, EDISON, NJ 08837 PH (732) 738-5454 FAX (732) 738-5841 EMAIL: 74201.3501@COMPUSERVE.COM

ANALYTICAL DATA REPORT

for

Lyondell Chemical Monaca, PA - 15061 Project: AHA / Monaca

Date Received: 7/23/99

Sample ID	Lab ID #
RC-ER-68-0799	R-6267.1
RC-ER-34-0799	R-6267.2
RC-EL-24-0799	R-6267.3
RC-EL-12-0799	R-6267.4
RC-EC-71-0799	R-6267.5
RC-EC-35-0799	R-6267.6
RC-EC-00-0799	R-6267.7
RC-EC-00-0799A	R-6267.8
Rinsate Blank	R-6267.9
Trip Blank	R-6267.10

These samples have been analyzed by EPA method 524.2 for a selected compound list. The results are not designed for use for drinking water purposes.

G. P. Kirpalani Manager

GPK/vb

R E L I A N C E LABORATORIES INC.



3090 WOODBRIDGE AVENUE, EDISON NJ 08837 PH (908) 738-5454 FAX (908) 738-5841

REDUCED LABORATORY DATA DELIVERABLES

Check if Complete

I.	Cover Page, Format, and Laboratory Certification (Include Cross Reference Table of Field I.D. and Lab I.D)	
II. III.	Chain of Custody Summary Sheets listing analytical results Including QA Day Information	ta V
IV.	Laboratory Chronicle and Methodology	
V.	Initial Calibration and Continuing Calibration	
VI.	Tune Summary (MS)	
VII.	Blank Summary	/
VIII.	Surrogate Recovery Summary	
IX.	Chromatograms / IR Spectra	
X.	Internal Standard Summary (MS)	
XI.	Matrix Spike / Spike Duplicate Summary	
XII.	Non-Conformance Summary	
	Laboratory Manager 7	14133 Date

Signature

RELIANCE LABORATORIES, INC.



175 MAY STREET, EDISON, NJ 08837 PH (732) 738-5454 FAX (732) 738-5841 EMAIL: 74201.3501@COMPUSERVE.COM

LABORATORY CHRONICLE

Customer Name Lyondell Chemical Date Received: 7/23/99 Date Sampled: 7/22/99 Sample ID: As per chain of custody
Organic Extraction:
1 Acids 2 Base / Neutrals 3 Pesticides/PCB's 4 TPHC
Analysis:
1 Volatiles 7/23/99 2 Acids 3 Base/Neutrals 4 Pesticides/PCB's 5 TPHC
1 Metals
2 Cyanides
Other Analysis:
Supervisor Review & Approval

RELIANCE LABORATORIES INC.



3090 WOODBRIDGE AVENUE, EDISON NJ 08837 PH (908) 738-5454 FAX (908) 738-5841

NON-CONFORMANCE SUMMARY

Reliance Labs received 10 water sample including blanks for BTEXS by method (EPA 524.2) from Lyondell Chemical on 23 July 1999. Samples consisted of 10 vials.

Matrix spike recovery analysis was performed on diluted sample since not enough sample was provided.

All analyses were performed within the required holding time.

STANDARD OPERATING PROCEDURE **METHOD 524.2**

1. Scope .

This is the general method for the procedure used to identify purgeable volatile organics in portable water. The sample is purged with ultra high purity helium and concentrated into a trap. The volatiles are then thermally desorbed onto a megabore column and identified using a mass spectrometer detector.

- **Equipment and Apparatus** 2.
- Sample containers- 40ml screw caps amber vials. A.
- Purge and Trap System. B.
 - 25cm VOCARB 3000 trap. 1.
- C. Glassware
 - 20 ml fritted purging vessels. 1.
 - 25 ml teflon sealed syringe with lever lock assembly. 2.
 - 10 μL syringes. 3.
- Gas Chromographic / Mass Spectrometer. D.
 - J&W Column type 1.

75 m, 0.53 mm ID, DB624 3 microns

- **Apparatus Conditions** E.
 - Tekmar (purge and trap) 1.
 - 2 min. Purge time a.
 - 250° for 2 min. Desorb time and temp.: Ъ.
 - 260° for 12 min. Bake time and temp. C.
 - 15 cc/min. d. Flow rate
 - 2. **GC** Conditions
 - 15 cc/min. Column flow a. 35° C
 - Initial temp. b.
 - 6° C/min. Ramping Rate C.
 - 200° C Final temp. d.
 - 47.25 min. Run time e.
 - Initial time 6 min.
- Stock Standards 3.
- Internal Standard A.
 - Flourobenzene 1.

f.

- B. Surrogates
 - 1. 1,2-dichlorobenzene-d4
 - 4-bromoflurobenzene
- Prepare standard solutions for all target compounds and surrogates at 20 ppm. C.
- Prepare internal standard at 20 ppm in methanol. D.
 - Prepare all standards and store in teflon sealed 1 ml vials.

4. Run Sequence

- A. Tune Instrument
- 1. Inject 1µL of 25 ppm BFB into GC.
 - a. Tune must pass against criteria.
 - b. Tune must be run before any samples, blank or calibrations can be run.
 - c. From time to tune 12 hours are available to run all QC data and samples.

B. Three Point Calibration Curve

- 1. Purge five (3) concentrations of standard solutions containing all the target analysis at 1 ppb, 2 ppb, 5 ppb.
- 2. The above standard must be run within 12 hours of injecting the BFB tune.
- 3. Created a calibration curve with the above standard runs.
 - a. If the 30% RSD deviation is exceeded the standards must be run again (still within 12 hours)
- 4. Create an identification file from this calibration curve for automated quantification.
- C. If time remains in the 12-hour run period go to step F.
- D. If the 12-hour period has expired, a new tune must be injected and a new sequence must be started.
- E. Once an initial calibration curve is established a continuing calibrations check may be run. A continuing calibration check is required every time the mass spectrometer is tuned.
 - 1. 2 ppb concentration of all target compounds is purged and quanted against current ID file.
 - 2. Check the response factors of this run against the average RF of the calibration file. The RF of the continuing calibration must be within \pm 50% D (difference) of the 5 point for all compounds.
 - The area counts of internal standard and surrogates must not be decreased by >30% from the most recent continuing calibration standard nor decrease by >50% from the initial calibration standard.

F. Daily Blank

- 1. Purge 20 ml of laboratory reagent water (nanopure) with 5 ppb internal standard and 5 ppb each surrogate.
- 2. Run this blank and quant against current ID file.
- 3. If blank does not meet criteria, it must be rerun before analyzing any samples.

G. Samples

- 1. Fill 25 ml syringe until it overflows with sample. Then adjust the volume to 20 ml exactly.
- 2. Inject 5 μ l each 25 ppm internal standard and surrogate standard solution into each sample.
- 3. Run and quant against the current 5 point calibration curves.
- 4. Any sample with target compound over 5 ppb must be rerun at the appropriate dilution
- 5. Any sample not injected in 12-hour period must be rerun.

H. Quality Control Sample (QCS)

1. Analyze a QCS from an external source at least quarterly.

R E L I A N C E LABORATORIES INC.



175 MAY STREET, EDISON, NJ 08837 PH (732) 738-5454 FAX (732) 738-5841 EMAIL: 74201.3501@COMPUSERVE.COM

<u>LABORATORY ID</u> NJ DEP NO. 12687 PA DER NO. 68437

CERTIFICATE OF ANALYSIS

Customer:

Lyondell Chemical

Sample:

Aqueous Samples

Date Sampled:

22 July 1999

Lab ID:

R-6267

Reference:

AHA / Monaca

26 July 1999

Units: μg/L

Sample ID	Benzene	Toluene	Ethylbenzene	Xylene	Styrene
	< 0.13	< 0.6	< 0.22	< 0.22	< 0.58
RC-ER-68-0799	< 0.13	< 0.6	< 0.22	< 0.22	< 0.58
RC-ER-34-0799	< 0.13	< 0.6	< 0.22	< 0.22	< 0.58
RC-EL-24-0799	< 0.13	< 0.6	< 0.22	< 0.22	< 0.58
RC-EL-12-0799	< 0.13	< 0.6	< 0.22	< 0.22	< 0.58
RC-EC-71-0799	< 0.13	< 0.6	< 0.22	< 0.22	< 0.58
RC-EC-35-0799		< 0.6	< 0.22	< 0.22	< 0.58
RC-EC-00-0799	< 0.13	< 0.6	< 0.22	< 0.22	< 0.58
RC-EC-00-0799A	< 0.13		< 0.22	< 0.22	< 0.58
Rinsate Blank	< 0.13	< 0.6	< 0.22	< 0.22	< 0.58
Trip Blank	< 0.13	< 0.6	< 0.22	~ U.ZZ	1 0.00

G. P. Kirpalani Manager

Data File : c:\hpchem\1\data\v6287.d

Acq On : 23 Jul 99 1:38 pm

Vial: 9 Operator: vb

Inst : 5971 - In

Sample : R-6267.1 Misc : AHA - RC-ER-68-0799

Multiplr: 1.00

Quant Time: Jul 26 9:08 1999

Method

: C:\HPCHEM\1\METHODS\RUN524.M

Title

: 524.2 Purgable Organics

Last Update : Mon Jul 26 09:07:31 1999

Response via: Multiple Level Calibration

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene	12.55	96	893237	5.00 ug/L	-0.01
System Monitoring Compounds 43) 4-bromofluorobenzene 55) 1,2-dichlorobenzene-d4	25.70 30.90	95 152	238029 148753	%R 5.09 ug/L 5.11 ug/L	
Target Compounds					Qvalue

^{(#) =} qualifier out of range (m) = manual integration Mon Jul 26 09:09:04 1999 v6287.d RUN524.M

Data File : c:\hpchem\1\data\v6287.d

: 23 Jul 99 1:38 pm Acq On

: R-6267.1

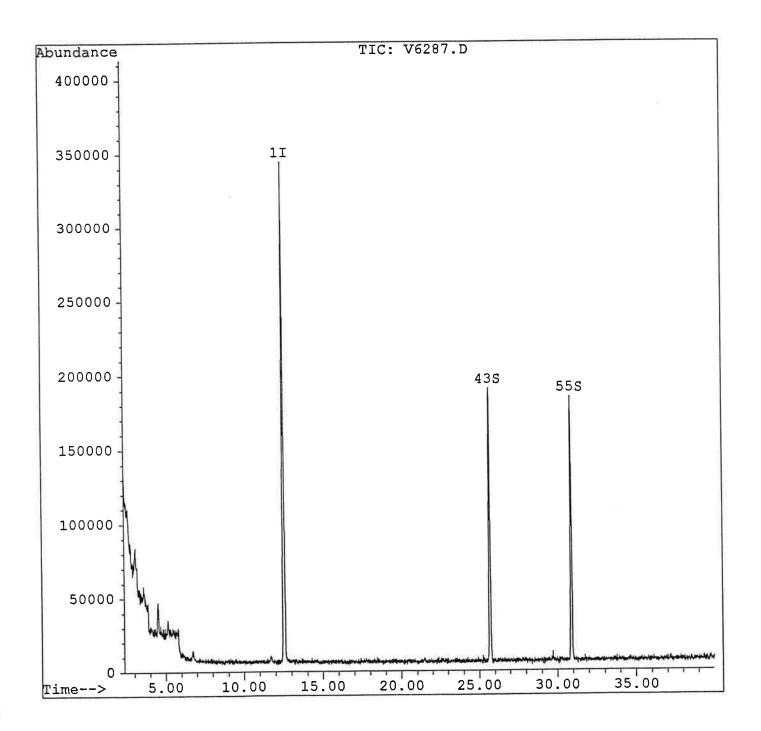
Sample : AHA - RC-ER-68-0799 Misc Quant Time: Jul 26 9:08 1999

Vial: 9 Operator: vb

: 5971 - In Inst

Multiplr: 1.00

: C:\HPCHEM\1\METHODS\RUN524.M Method



Vial: 10

Inst : 5971 - In

Operator: vb

Multiplr: 1.00

Data File : c:\hpchem\1\data\v6288.d

: 23 Jul 99 2:23 pm Acq On

: R-6267.2 Sample

Method

Misc : AHA - RC-ER-34-0799

Quant Time: Jul 26 9:09 1999

: C:\HPCHEM\1\METHODS\RUN524.M

Title : 524.2 Purgable Organics
Last Update : Mon Jul 26 09:07:31 1999

Response via : Multiple Level Calibration

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene	12.55	96	856616	5.00 ug/L	0.00
System Monitoring Compounds 43) 4-bromofluorobenzene 55) 1,2-dichlorobenzene-d4	25.69 30.90	95 152	228539 143444	%F 5.10 ug/L 5.14 ug/L	
Target Compounds					Qvalue

^{(#) =} qualifier out of range (m) = manual integration v6288.d RUN524.M Mon Jul 26 09:09:27 1999

Vial: 10

Inst : 5971 - In

Operator: vb

Multiplr: 1.00

Data File : c:\hpchem\1\data\v6288.d

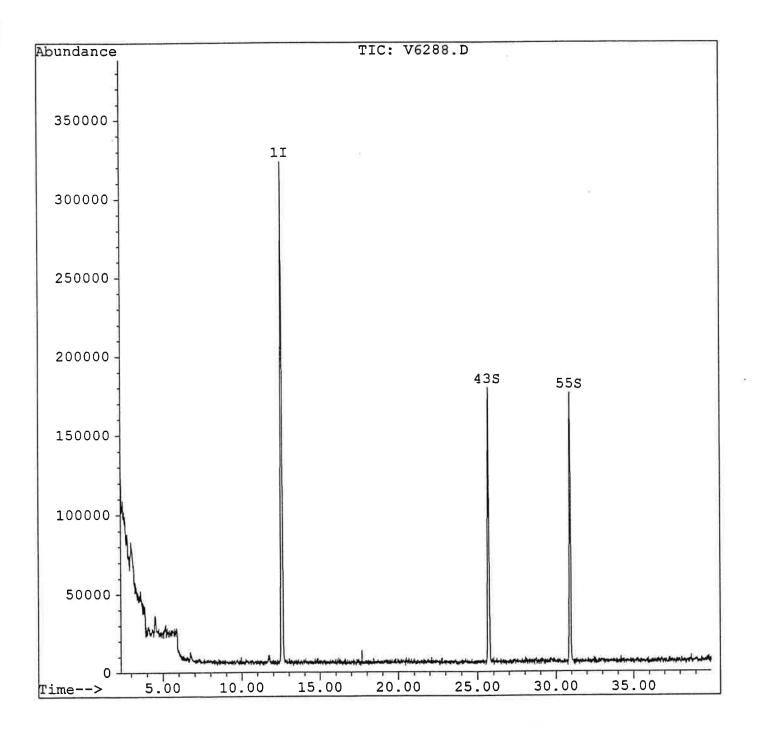
: 23 Jul 99 2:23 pm Acq On

: R-6267.2 Sample

: AHA - RC-ER-34-0799 Misc

Quant Time: Jul 26 9:09 1999

: C:\HPCHEM\1\METHODS\RUN524.M Method



Data File : c:\hpchem\1\data\v6289.d

Acq On : 23 Jul 99 3:10 pm

Vial: 11

Operator: vb

Sample : R-6267.3 Misc : AHA - RC-EL-24-0799 : R-6267.3

Inst : 5971 - In Multiplr: 1.00

Quant Time: Jul 26 9:09 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M
Title : 524.2 Purgable Or

Last Update : Mon Jul 26 09:07:31 1999

Response via : Multiple Level Calibration

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene	12.56	96	885470	5.00 ug/L	0.00
System Monitoring Compounds 43) 4-bromofluorobenzene 55) 1,2-dichlorobenzene-d4	25.69 30.90	95 152	234115 145315	%R 5.05 ug/L 5.04 ug/L	
Target Compounds					Qvalue

^(#) = qualifier out of range (m) = manual integration Mon Jul 26 09:10:03 1999 v6289.d RUN524.M

Vial: 11

: 5971 - In

Operator: vb

Multiplr: 1.00

Inst

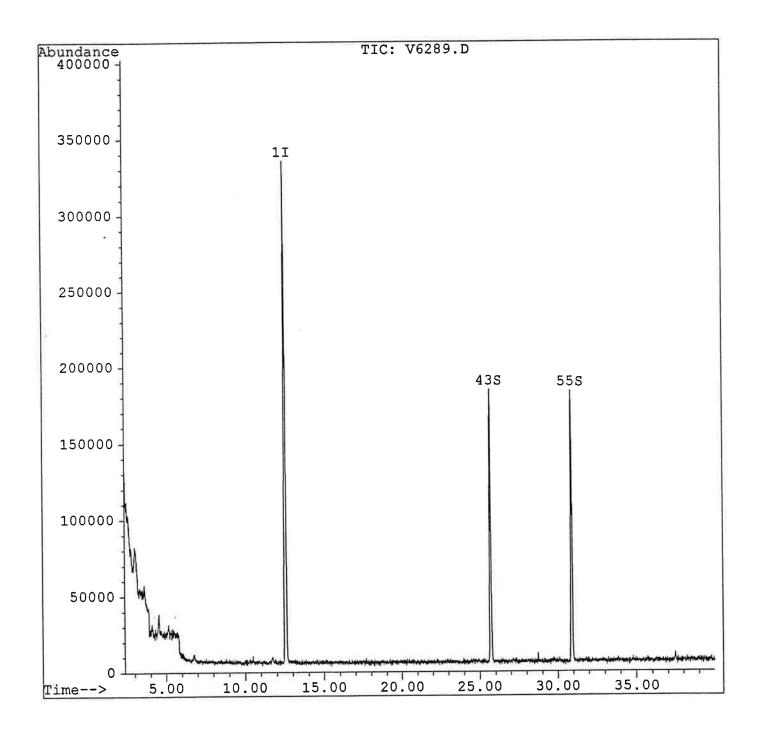
Data File : c:\hpchem\1\data\v6289.d

Acq On : 23 Jul 99 3:10 pm

Sample : R-6267.3

Misc : AHA - RC-EL-24-0799 Quant Time: Jul 26 9:09 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M



Data File : c:\hpchem\1\data\v6290.d

Acq On : 23 Jul 99 3:55 pm

Vial: 12

Operator: vb

Inst : 5971 - In Multiplr: 1.00

Sample : R-6267.4 Misc : AHA - RC-: AHA - RC-EL-12-0799

Quant Time: Jul 26 9:10 1999

: C:\HPCHEM\1\METHODS\RUN524.M

Method Title : 524.2 Purgable Organics Last Update : Mon Jul 26 09:07:31 1999 Response via : Multiple Level Calibration

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene	12.54	96	849200	5.00 ug/L	-0.02
System Monitoring Compounds 43) 4-bromofluorobenzene 55) 1,2-dichlorobenzene-d4	25.70 30.91	95 152	234296 145039	%R 5.27 ug/L 5.24 ug/L	
Manual Compounds					Ovalue

Target Compounds

^{(#) =} qualifier out of range (m) = manual integration Mon Jul 26 09:10:31 1999 v6290.d RUN524.M

Vial: 12 Operator: vb

Multiplr: 1.00

Inst : 5971 - In

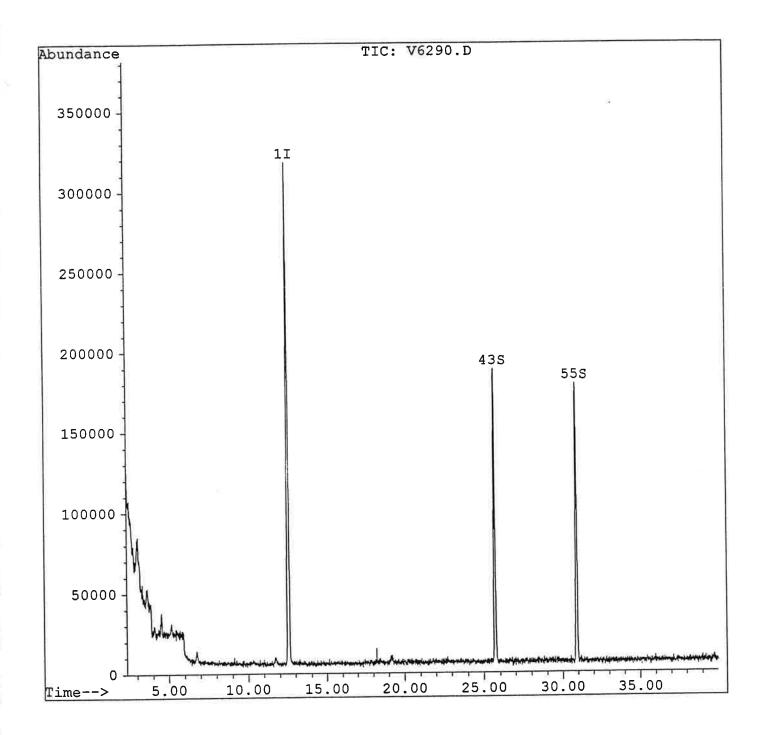
Data File : c:\hpchem\1\data\v6290.d

3:55 pm : 23 Jul 99 Acq On

: R-6267.4 Sample

: AHA - RC-EL-12-0799 Misc Quant Time: Jul 26 9:10 1999

: C:\HPCHEM\1\METHODS\RUN524.M Method



Data File : c:\hpchem\1\data\v6291.d

Acq On : 23 Jul 99 4:42 pm

Vial: 13

Operator: vb

Sample : R-6267.5 Misc : AHA - RC-EC-71-0799

Inst : 5971 - In Multiplr: 1.00

Quant Time: Jul 26 9:10 1999

: C:\HPCHEM\1\METHODS\RUN524.M

Title

: 524.2 Purgable Organics

Last Update : Mon Jul 26 09:07:31 1999

Response via : Multiple Level Calibration

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene	12.56	96	839688	5.00 ug/L	0.00
System Monitoring Compounds 43) 4-bromofluorobenzene 55) 1,2-dichlorobenzene-d4	25.69 30.92	95 152	226420 141637	%R 5.15 ug/L 5.18 ug/L	
Target Compounds					Qvalue

^{(#) =} qualifier out of range (m) = manual integration v6291.d RUN524.M Mon Jul 26 09:10:54 1999

Data File : c:\hpchem\1\data\v6291.d

Acq On : 23 Jul 99 4:42 pm

Sample : R-6267.5

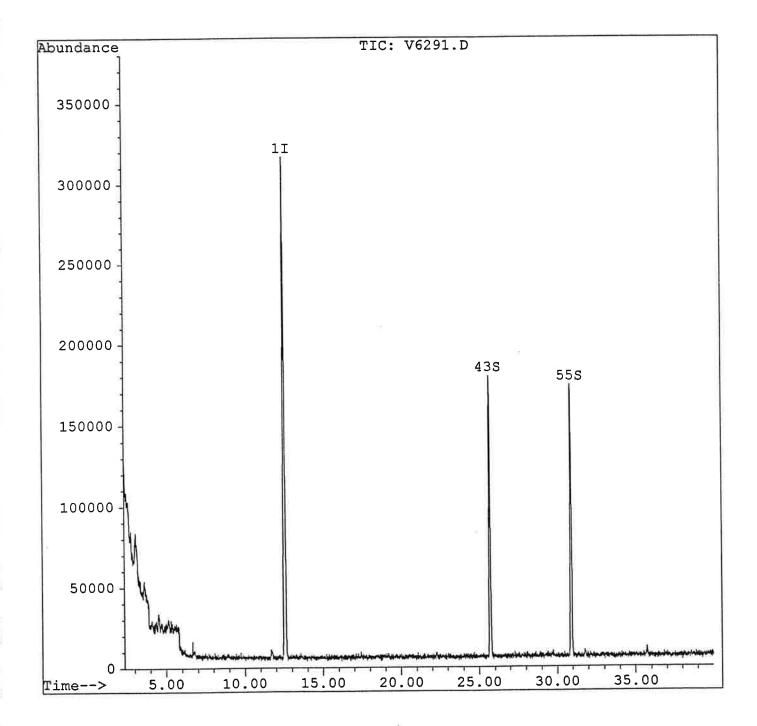
Misc : AHA - RC-EC-71-0799 Quant Time: Jul 26 9:10 1999 Vial: 13
Operator: vb

Inst : 5971 - In

Multiplr: 1.00

Method : C:\HPCHEM\1\METHODS\RUN524.M
Title : 524.2 Purgable Organics

Last Update : Mon Jul 26 09:07:31 1999 Response via : Multiple Level Calibration



Data File : c:\hpchem\1\data\v6292.d

: 23 Jul 99 5:27 pm

Vial: 14 Operator: vb

Acq On

Inst : 5971 - In

: R-6267.6 Sample

Misc : AHA - RC-EC-35-0799

Multiplr: 1.00

Quant Time: Jul 26 9:11 1999

: C:\HPCHEM\1\METHODS\RUN524.M Method

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene	12.54	96	905814	5.00 ug/L	-0.02
System Monitoring Compounds 43) 4-bromofluorobenzene 55) 1,2-dichlorobenzene-d4	25.68 30.90	95 152	244932 150055	%R 5.17 ug/L 5.08 ug/L	Recovery 103.31% 101.69%
Target Compounds					Qvalue

^(#) = qualifier out of range (m) = manual integration v6292.d RUN524.M Mon Jul 26 09:11:19 1999

Vial: 14

: 5971 - In

Operator: vb

Inst

Data File : c:\hpchem\1\data\v6292.d

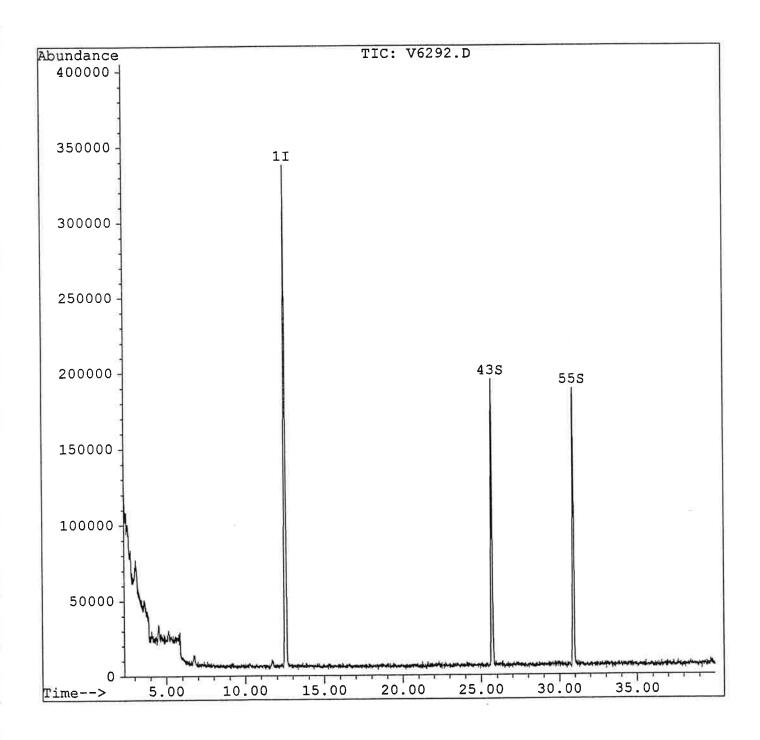
: 23 Jul 99 5:27 pm Acq On

: R-6267.6 Sample

Misc Quant Time: Jul 26 9:11 1999

Multiplr: 1.00 : AHA - RC-EC-35-0799

: C:\HPCHEM\1\METHODS\RUN524.M Method



Data File : c:\hpchem\1\data\v6293.d

Acq On : 23 Jul 99 6:14 pm

Vial: 15 Operator: vb

Inst : 5971 - In Multiplr: 1.00

Sample : R-6267.7 Misc : AHA - RC-EC-00-0799

Quant Time: Jul 26 9:11 1999

: C:\HPCHEM\1\METHODS\RUN524.M Method

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene	12.55	96	848054	5.00 ug/L	0.00
System Monitoring Compounds 43) 4-bromofluorobenzene 55) 1,2-dichlorobenzene-d4	25.69 30.90	95 152	224573 136328	5.06 ug/L	
Target Compounds					Qvalue

^{(#) =} qualifier out of range (m) = manual integration Mon Jul 26 09:11:41 1999 v6293.d RUN524.M

Data File : c:\hpchem\1\data\v6293.d

Acq On : 23 Jul 99 6:14 pm

: R-6267.7 Sample

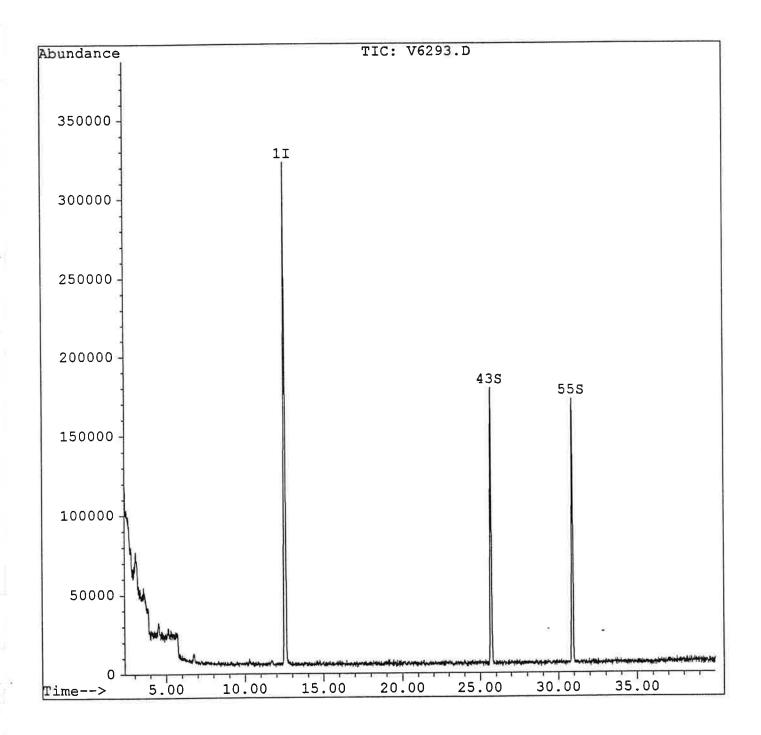
: AHA - RC-EC-00-0799 Misc Quant Time: Jul 26 9:11 1999

Operator: vb : 5971 - In Inst

Multiplr: 1.00

Vial: 15

: C:\HPCHEM\1\METHODS\RUN524.M Method



Data File : c:\hpchem\1\data\v6294.d

Acq On : 23 Jul 99 7:01 pm

Vial: 16 Operator: vb Inst : 5971 - In Multiplr: 1.00

: R-6267.8 Sample

Misc : AHA - RC-EC-00-0799A

Quant Time: Jul 26 9:11 1999

: C:\HPCHEM\1\METHODS\RUN524.M Method

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene	12.55	96	839260	5.00 ug/L	0.00
System Monitoring Compounds 43) 4-bromofluorobenzene 55) 1,2-dichlorobenzene-d4	25.69 30.90	95 152	227812 139566	%R 5.19 ug/L 5.10 ug/L	
Target Compounds					Qvalue

^(#) = qualifier out of range (m) = manual integration v6294.d RUN524.M Mon Jul 26 09:11:59 1999

Vial: 16

: 5971 - In

Operator: vb

Multiplr: 1.00

Inst

Data File : c:\hpchem\1\data\v6294.d

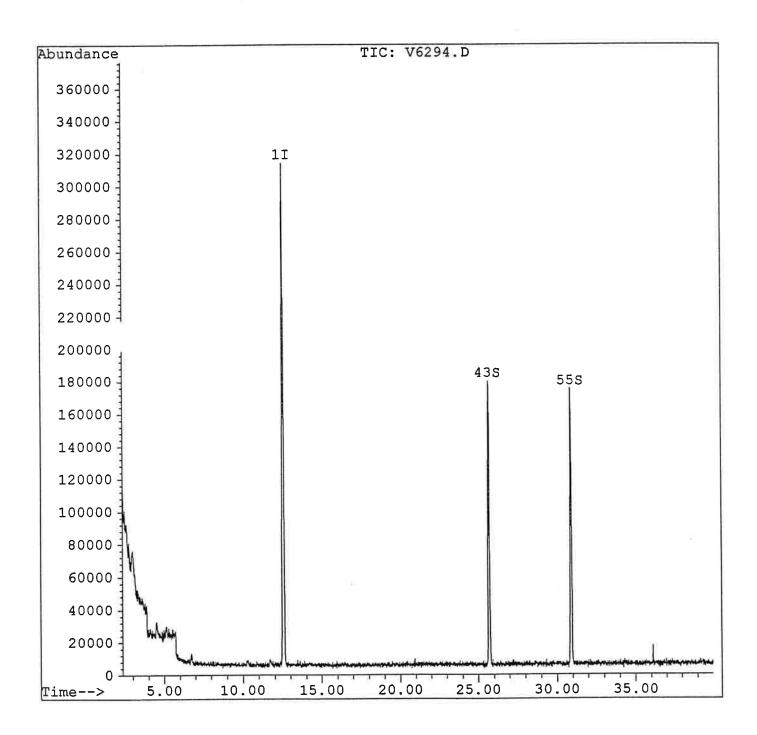
: 23 Jul 99 7:01 pm Acq On

: R-6267.8 Sample

: AHA - RC-EC-00-0799A Misc

Quant Time: Jul 26 9:11 1999

: C:\HPCHEM\1\METHODS\RUN524.M Method



Vial: 4

Multiplr: 1.00

: 5971 - In

Operator: vb

Inst

Data File : C:\HPCHEM\1\DATA\V6296.D

: 23 Jul 99 8:34 pm Acq On

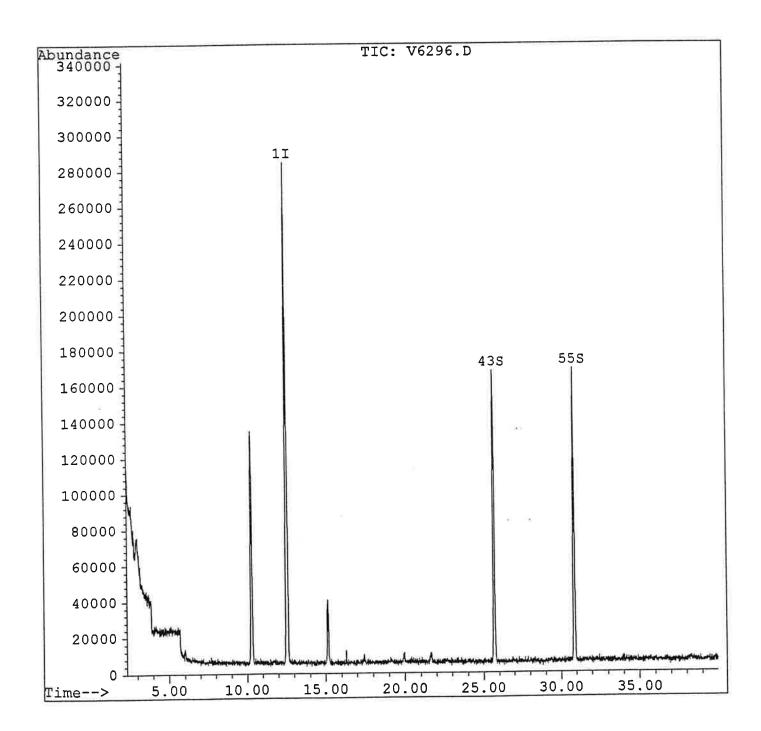
: R-6267.10 Sample

: AHA - Trip Blank Misc

Quant Time: Jul 26 10:14 1999

: C:\HPCHEM\1\METHODS\RUN524.M

Method : 524.2 Purgable Organics Title Last Update : Mon Jul 26 09:07:31 1999 Response via : Multiple Level Calibration



Data File : C:\HPCHEM\1\DATA\V6295.D

Acq On : 23 Jul 99 7:47 pm

Vial: 3

Operator: vb Inst : 5971 - In

Multiplr: 1.00

Sample : R-6267.9 Misc : AHA - Rinsate Blank

Quant Time: Jul 26 10:15 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M
Title : 524.2 Purgable Organics
Last Update : Mon Jul 26 09:07:31 1999 Response via : Multiple Level Calibration

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene	12.54	96	805721	5.00 ug/L	-0.02
System Monitoring Compounds 43) 4-bromofluorobenzene 55) 1,2-dichlorobenzene-d4	25.69 30.90	95 152	227201 142647	%R 5.39 ug/L 5.43 ug/L	
Target Compounds					Qvalue

^{(#) =} qualifier out of range (m) = manual integration V6295.D RUN524.M Mon Jul 26 10:15:16 1999

WATER VOLATILE SYSTEM MONITORING COMPOUND RECOVERY

Customer : Lyondell

Γ		SMC1	SMC2		OTHER	TOT
- 1	SAMPLE NO.	#	#	#	#	OUT
01	VBLK01	97	100			
02	R-6267.1	102	102			
03	R-6267.2	102	103			
04	R-6267.3	101	101			
05	R-6267.4	105	105			
06	R-6267.5	103	104			
07	R-6267.6	103	102			
80	R-6267.7	101	99			
09	R-6267.8	104	102			
10	R-6267.9	108	112			
11	R-6267.10	109	109			
12	VBLK02	118	122 *			
13	R-6267.1MS	98	100			
14	R-6267.1MSD	97	96			
15						
16						
17						
18						
19						
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21						
22						
23						
24						
25						
26						
27						
28				1		
29						
30						

QC LIMITS (80-120)

SMC1 = 4-Bromofluorobenzene

SMC2 = 1,2-dichlorobenzene-d4

(80-120)

- # Column to be used to flag recovery values
- * Values outside of contract required QC limits
- D System Monitoring Compound diluted out

FORM II VOA-1

WATER SEMIVOLATILE MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

R-6267.1 Matrix Spike - Sample No.:

COMPOUND	SPIKE ADDED (ug/Kg)	SAMPLE CONCENTRATION (ug/Kg)	MS CONCENTRATION (ug/Kg)	MS % REC #	
Benzene	3.00	0.00	2.89	96	(80-120)
Toluene	3.00	0.00	2.82	94	(80-120)
Ethylbenzene	3.00	0.00	2.95	98	(80-120)
m&p-xylenes	3.00	0.00	2.91	97	(80-120)
o-xylenes	3.00	0.00	2.93	98	(80-120)
Styrene	3.00	0.00	2.90	97	(80-120)

	SPIKE ADDED	MSD CONCENTRATION	MSD %	%	QC L	IMITS
COMPOUND	(ug/Kg)	(ug/Kg)	REC #	RPD #	RPD	REC.
Benzene	3.00	2.88	96	0	20	(80-120)
Toluene	3.00	2.79	93	1	20	(80-120)
Ethylbenzene	3.00	2.92	97	1	20	(80-120)
m&p-xylenes	3.00	2.93	98	1	20	(80-120)
o-xylenes	3.00	2.89	96	1	20	(80-120)
Styrene	3.00	2.88	96	1	20	(80-120)

Column to be used to flag recovery and RPD values with an asterisk * Values outside of QC limits

Comments:	

VOLATILE METHOD BLANK SUMMARY

VBLK0

	Customer :	Lyondell	-		
Lab File ID V62	86.D			Lab Sample ID:	BLANK1
Date Analyzed:	7/23/99			Time Analyzed:	1251
GC Column:		0.53 (mm)		
			•		
Instrument ID:				04110150 110 1	ND MOD.
THI	S METHOD BLANK	APPLIES TO THE	FOLLOWING	SAMPLES, MS A	ND M2D:
		LAB	LAB	TIME	
	SAMPLE NO.	SAMPLE ID	FILE ID	ANALYZED	
	R-6267.1	ER-68	V6287.D	1338	
	R-6267.2	ER-34	V6288.D	1423	
	R-6267.3	EL-24	V6289.D	1510 1555	
	R-6267.4	EL-12	V6290.D V6291.D	1642	
	R-6267.5	EC-71 EC-35	V6291.D	1727	
	R-6267.6	EC-00	V6293.D	1814	
	R-6267.7 R-6267.8	EC-00A	V6294.D	1901	
	R-6267.9	RINSATE	V6295.D	1947	
	R-6267.10	TBLANK	V6296.D	2034	
11					
12					
13					
14					
15					
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24					
25 26			+		5
27					
28					
29					
30					
COMMENTS:					

Page 1 of 1

FORM IV VOA

Data File : c:\hpchem\1\data\v6286.d

Acq On : 23 Jul 99 12:51 pm

Vial: 8 Operator: vb

Inst : 5971 - In

Sample : blank
Misc : blank

Multiplr: 1.00

Quant Time: Jul 26 9:08 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M
Title : 524.2 Purgable Organics

Last Update : Mon Jul 26 09:07:31 1999

Response via : Multiple Level Calibration

Internal Standards	R.T.	QIon	Response	Conc Units Dev(Min)
1) Fluorobenzene	12.56	96	890100	5.00 ug/L 0.00
System Monitoring Compounds 43) 4-bromofluorobenzene 55) 1,2-dichlorobenzene-d4	25.68 30.90	95 152	225933 145114	%Recovery 4.85 ug/L 96.98% 5.00 ug/L 100.08%
Target Compounds				Qvalue

^{(#) =} qualifier out of range (m) = manual integration Mon Jul 26 09:08:42 1999 v6286.d RUN524.M

Vial: 8

Multiplr: 1.00

: 5971 - In

Operator: vb

Inst

Data File : c:\hpchem\1\data\v6286.d

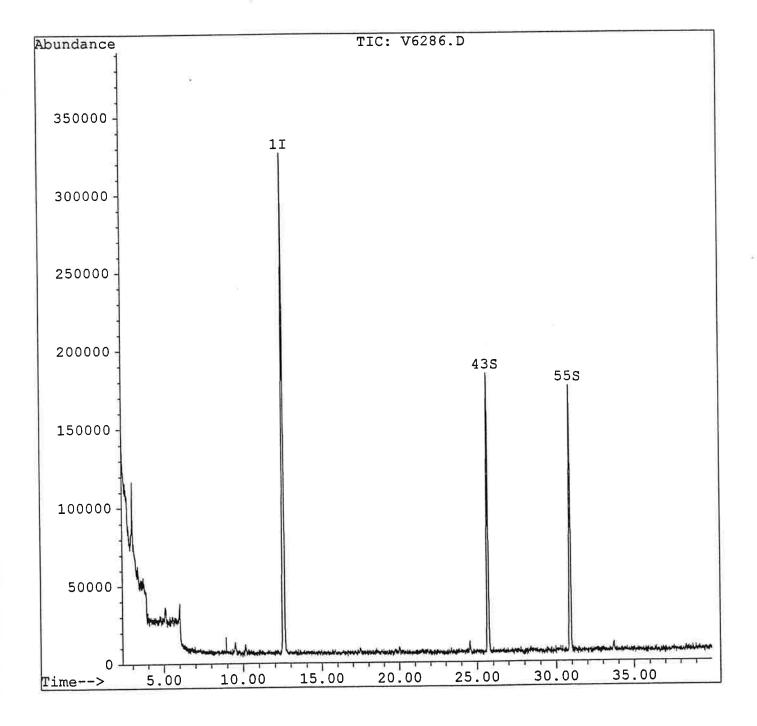
: 23 Jul 99 12:51 pm

Sample : blank : blank Misc

Quant Time: Jul 26 9:08 1999

: C:\HPCHEM\1\METHODS\RUN524.M Method

: 524.2 Purgable Organics Title Last Update : Mon Jul 26 09:07:31 1999 Response via : Multiple Level Calibration



VOLATILE METHOD BLANK SUMMARY

VBLK0

	Customer :	Lyondell	-				
Lab File ID V62	99.D			Lab Sample ID	: BLANK2		
Date Analyzed:	7/27/99			Time Analyzed	: 1319		
GC Column:	DB-624 ID:	0.53 (mm))				
Instrument ID:	HP5971						
	THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:						
		LAB	LAB	TIME			
	SAMPLE NO.	SAMPLE ID	FILE ID	ANALYZED			
	R-6267.1MS	ER-68MS	V6300.D	1406			
	R-6267.1MSD	ER-68MSD	V6302.D	1620			
03							
04							
05							
06							
07 08							
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30			1		l		
COMMENTS:							
(

Page 1 of 1

J31

FORM IV VOA

Data File : C:\HPCHEM\1\DATA\V6299.D

Acq On : 27 Jul 99 1:19 pm Sample : blank Misc : blank

Vial: 3 Operator: vb

Inst : 5971 - In

Multiplr: 1.00

Quant Time: Jul 27 15:54 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M
Title : 524.2 Purgable Organics

Last Update : Tue Jul 27 15:54:47 1999

Response via : Multiple Level Calibration

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene	12.55	96	690549	5.00 ug/L	0.00
System Monitoring Compounds 43) 4-bromofluorobenzene 55) 1,2-dichlorobenzene-d4	25.71 30.91	95 152	212666 137186	%R 5.88 ug/L 6.10 ug/L	
Target Compounds					Qvalue

^{(#) =} qualifier out of range (m) = manual integration Tue Jul 27 15:55:00 1999 V6299.D RUN524.M

Data File : C:\HPCHEM\1\DATA\V6299.D

1:19 pm : 27 Jul 99 Acq On

Vial: 3 Operator: vb Inst : 5971 - In

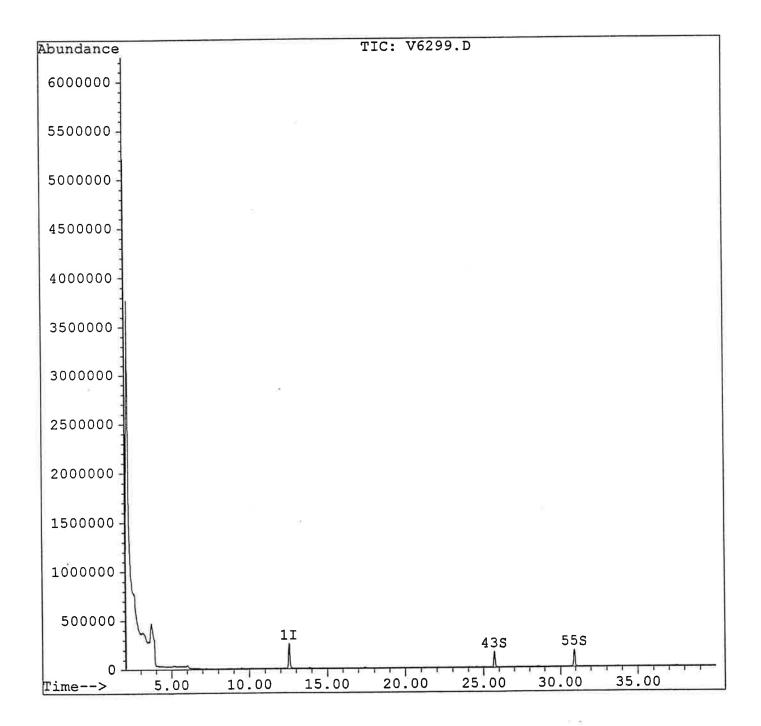
: blank Sample : blank Misc

Multiplr: 1.00

Quant Time: Jul 27 15:54 1999

: C:\HPCHEM\1\METHODS\RUN524.M

: 524.2 Purgable Organics Title Last Update : Tue Jul 27 15:54:47 1999 Response via : Multiple Level Calibration



VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK BROMOFLUOROBENZENE (BFB)

Customer:	Lyondell	
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Lab File ID: V6282.D

BFB Injection Date: 7/23/99

Instrument ID: HP5971A

BFB Injection Time: 0845

GC Column: DB-624

ID: 0.53 (mm)

		%RELATIVE		
m/e	ION ABUNDANCE CRITERIA	ABUNDANC	Ξ	
50	8.0 - 40.0% of mass 95	19.9		
75	30.0 - 66.0% of mass 95	48.4		
95	Base peak, 100% relative abundance	100.0		
96	5.0 - 9.0% of mass 95	6.5		
173	Less than 2.0% of mass 174	0.0 (0.0)1	
174	50.0 - 120.0% of mass 95	63.2		
175	4.0 - 9.0% of mass 174	4.7 (7.4)1	
176	93.0 - 101.0% of mass 174	60.3 (95.5)1	
177	5.0 - 9.0% of mass 176	4.0 (6.6)2	

1-Value is % mass 174

2-Value is % mass 176

This check applies to the following SAMPLES, MS, MSD, BLANKS and STANDARDS:

SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
		V6283.D	7/23/99	1010
01 VSTD050	ICC002	V6284.D	7/23/99	1058
02 VSTD005	ICC005		7/23/99	1144
03 VSTD001	ICC001	V6285.D		
04 VBLK01	BLANK1	V6286.D	7/23/99	1251
05 R-6267.1	ER-68	V6287.D	7/23/99	1338
06 R-6267.2	ER-34	V6288.D	7/23/99	1423
07 R-6267.3	EL-24	V6289.D	7/23/99	1510
08 R-6267.4	EL-12	V6290.D	7/23/99	1555
09 R-6267.5	EC-71	V6291.D	7/23/99	1642
10 R-6267.6	EC-35	V6292.D	7/23/99	1727
11 R-6267.7	EC-00	V6293.D	7/23/99	1814
12 R-6267.8	EC-00A	V6294.D	7/23/99	1901
13 R-6267.9	RINSATE	V6295.D	7/23/99	1947
14 R-6267.10	TBLANK	V6296.D	7/23/99	2034
15				
16				
17				
18				
19				
20				
21				
22	-			

Page 1 of 1

FORM V VOA

Data File : C:\HPCHEM\1\DATA\V6282.D

Acq On : 23 Jul 99 8:45 am

Sample : bfb

Misc :

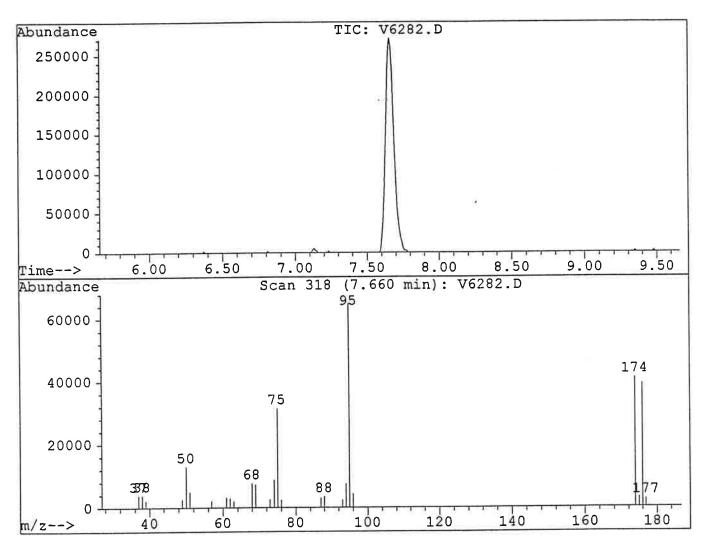
Vial: 1
Operator: vb

Inst : 5971 - In

Multiplr: 1.00

Method : C:\HPCHEM\1\METHODS\RUN524.M

Title : 524.2 Purgable Organics



Peak Apex is scan: 318

1	Target Mass	1	Rel. to Mass	 	Lower Limit%	1	Upper Limit%	1	Rel. Abn%	! 	Raw Abn	1	Result Pass/Fail	
1	50	1	95	ī	- 	1	40	1	19.9	1	12924	1	PASS	1
i	75	i	95	ıi.	3.0	i	80	î	48.4	ĺ	31392	1	PASS	- 1
i	95	i.	95	î	100	i	100	Î	100.0	ĺ	64920	1	PASS	- 1
i	96	i	95	i	5	i	9	î	6.5	Ī	4226	1	PASS	-1
1	173	i	174	i	Ō	i	2	î	0.0	ì	0	1	PASS	1
î	174	i	95	i	50	i	100	î	63.2	i	41000	1	PASS	1
Ť	175	Ť	174	-i	5	i	9	i	7.4	Í	3025	1	PASS	1
Ť	176	÷	174	i	95	i	101	i	95.5	i	39144	1	PASS	Î
i	177	i	176	ì	5	į	9	į	6.6	Ì	2575	I	PASS	l

VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK BROMOFLUOROBENZENE (BFB)

Customer : Lyondell	
Lab File ID: V6297.D	BFB Injection Date: 7/27/99
Instrument ID: HP5971A	BFB Injection Time: 1107
GC Column:DB-624	

		%RELATIVE				
m/e	ION ABUNDANCE CRITERIA	ABUNDANCE				
50	8.0 - 40.0% of mass 95	21.4				
75	30.0 - 66.0% of mass 95	43.9				
95	Base peak, 100% relative abundance	100.0				
96	5.0 - 9.0% of mass 95	6.2				
173	Less than 2.0% of mass 174	0.0 (0.0)1				
174	50.0 - 120.0% of mass 95	60.3				
175	4.0 - 9.0% of mass 174	4.2 (6.9)1				
176	93.0 - 101.0% of mass 174	59.3 (98.4)1				
177	5.0 - 9.0% of mass 176	4.2 (7.2)2				

177 | 5.0 - 9.0% of mass 176 1-Value is % mass 174

2-Value is % mass 176

This check applies to the following SAMPLES, MS, MSD, BLANKS and STANDARDS:

	LAB	LAB	DATE	TIME ANALYZED
SAMPLE NO.	SAMPLE ID	FILE ID	ANALYZED	
01 VSTD050	CC0022	V6298.D	7/27/99	1225
02 VBLK02	BLANK2	V6299.D	7/27/99	1319
03 R-6267.1MS	ER-68MS	V6300.D	7/27/99	1406
04 R-6267.1MSD	ER-68MSD	V6302.D	7/27/99	1620
05				V
06				
07				
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14		-		
15	 			
		-		
16				
17				
18				
19	-			
20				
21				
22				

Page 1 of 1

FORM V VOA

Data File : C:\HPCHEM\1\DATA\V6297.D

Acq On : 27 Jul 99 11:07 am

Sample : bfb

Misc :

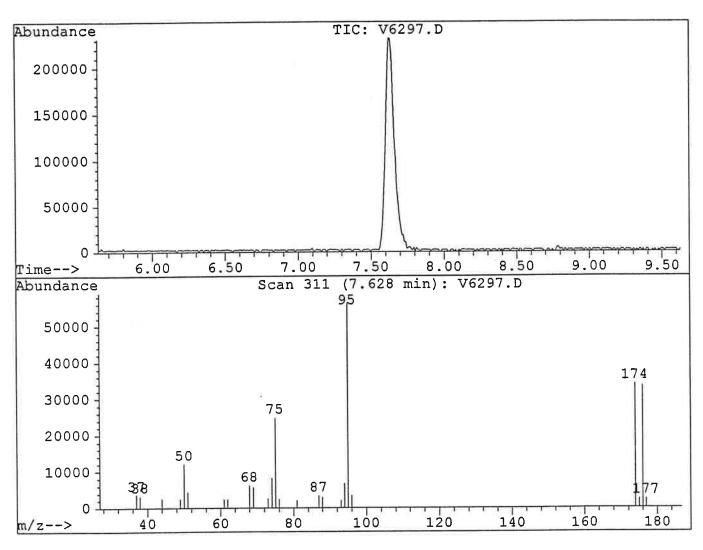
Vial: 1
Operator: vb

Inst : 5971 - In

Multiplr: 1.00

Method : C:\HPCHEM\1\METHODS\RUN524.M

Title : 524.2 Purgable Organics



Peak Apex is scan: 311

1	Target Mass		Rel. to Mass		Lower Limit%		Upper Limit%	1	Rel. Abn%		Raw Abn		Result Pass/Fail	
ī	50	ī	95	1	15	1	40	1	21.4	1	12065	1	PASS	1
i	75	î	95	Ĺ	30	1	80	1	43.9	1	24752	1	PASS	1
-i	95	i.	95	ì	100	Ì	100	ĺ	100.0	- 1	56368	1	PASS	1
i	96	ï	95	Î	5	ì	9	1	6.2	1	3468	1	PASS	1
i	173	i	174	Î	0	ì	2	Ĩ	0.0	- 1	0	1	PASS	1
i	174	i	95	i	50	Î	100	Î	60.3	Î	33976	1	PASS	1
i	175	ì	174	i	5	İ	9	Ì	6.9	Ĩ	2351	1	PASS	1
ì	176	i	174	i	95	Î	101	Ì	98.4	ĺ	33424	1	PASS	-1
Ì	177	i	176	i	5	i	9	İ	7.2	İ	2393	1	PASS	1

VOLATILE ORGANICS INITIAL CALIBRATION DATA

		Customer	Lyondell			
instrument ID:	HP5971A	_	Calibration Date(s):	7/23/99	7/23/99	
		= *:	Calibration Times:	1010	1144	

GC Column: DB-624	ID:	0.53_	(mm)			
Lab File ID: RRF05 = V6284.D	RRF01 =	V6285.D	j	RRF02 = V	6283.D	
COMPOUND	RRF01	RRF02	RRF05		RRF	% RSD
Benzene	1.359	1.217	1.141		1.239	8.9
Toluene	1.338	1.085	0.987		1.137	15.9
Ethylbenzene	1.288	1.134	1.075		1.166	9.4
m&p-xylenes	0.893	0.785	0.750		0.809	9.2
o-xylene	0.934	0.790	0.746		0.823	11.9
Styrene	0.649	0.560	0.540		0.583	10.0
4-bromofluorobenzene	0.265	0.255	0.265		0.262	2.2
1,2-dichlorobenzene-d4	0.166	0.160	0.162		0.163	1.9

Page 1 of 1

FORM VI VOA

VOLATILE CONTINUING CALIBRATION CHECK

Customer : Lyondell

Instrument ID: HP5971A Calibration Date: 7/27/99

Time: 1225

Lab File ID: V6298.D Init. Calib. Date(s): 7/23/99 7/23/99

Init. Calib. Times: 1010 1144

GC Column:

DB-624

ID: 0.53 (mm)

			MIN	~ 5	MAX %D
COMPOUND	RRF	RRF20	RRF	%D	%D
Benzene	1.239	1.191		3.9	
Toluene	1.137	1.107		2.6	
Ethylbenzene	1.166	1.173		-0.6	
m&p-xylenes	0.809	0.818		-1.1	ļ
o-xylene	0.823	0.804		2.3	
Styrene	0.583	0.564		3.3	-
					-
					-
					-
		-			
				ļ	+
					-
12					
					-
4-Bromofluorobenzene	0.262	0.252	-	3.8	-
1,2-dichlorobenzene-d4	0.163	0.157	1	3.5	

All other compounds must meet a minimum RRF of 0.010.

VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Customer : Lyondell

Lab File ID (Standard): V6298.D

Date Analyzed: 7/27/99

Instrument ID: HP5971A

Time Analyzed: 1225

GC Column: DB-624

ID: 0.53 (mm)

								-	-	
	IS1	DT	ADEA	#	RT	#	AREA	#	RT	#
	AREA #	RT #	AREA	#	KI_	#	AREA		IXI	
12 HOUR STD	878256	12.55		_						
UPPER LIMIT	1756512	13.05		_						
LOWER LIMIT	439128	12.05		_						
SAMPLE										
NO.										
01 VBLK02	690549	12.55								
02 R-6267.1MS	857144	12.55								
03 R-6267.1MSD	865445	12.57								
04										
05										
06										
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08										
09								-		
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12									-	
13			2							
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15				_					-	
16								_		_
17						_				
18										
19								-		
20										
21										
22							L			

IS1 = Fluorobenzene

AREA UPPER LIMIT = +100% of internal standard area AREA LOWER LIMIT = -50% of internal standard area RT UPPER LIMIT = +0.50 minutes of internal standard RT RT LOWER LIMIT = -0.50 minutes of internal standard RT

Column used to flag values outside QC limits with an asterisk

* Values outside of QC limits.

Page 1 of 1

FORM VIII VOA

DEPARTMENT OF ENVIRONMENTAL PROTECTION

Certifies That
Rellance Laboratories, Inc.
3090 Wood Bridge Avenue
Edison, NJ 08837



ויטו ניות

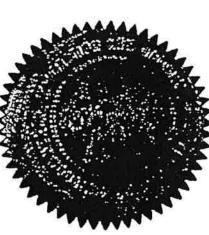
having duly met the requirements of the

Regulations Governing Laboratory Certification And Standards Of Performance NJ.A.C. 7:18 et. seq.

is hereby approved as a

State Certified Water Laboratory

To perform the analyses as indicated on the Annual Certified Parameter List which must accompany this certificate to be valid



DEPARTMENT OF ENVIRONMENTAL PROTECTION

12687
PERMANENT CERTIFICATION NUMBER

. Ianiiary 11 1989
DATE

N.J.A.C. 7:18-2.11(d) and agreed to by the Laboratory Manager on filing the application This certification is subject to unannounced laboratory inspections as specified by

TO BE CONSPICUOUSLY DISPLAYED AT THE LABORATORY WITH THE ANNUAL CERTIFIED PARAMETER LIST.

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400 Frankfort Rd, Monaca, Pa 15061 Applied

400 Frankfort Rd Monaca Pa., 15061

(724) 728 - 6586 (724) 728 - 6498

> Phone **Fa**x:

Lyondell Chemical

Customer: Address: b_petroff@worldnet.att.net

E-Mail:

Associates, Inc. Hydrology

Page 1 of 1 Trun around time Standard (stanrard / rush) Date: 7/22/197 LAB ID: 8-6267 Project ID:

Fax results: Y N E-Mail results: Y N

Sampler Name: BDP Preserved Y N Sample Intact: Y N

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		Dissolved Metals	
1	METALS	Total Metals (list below)	
	ME	Priority Pol. (13)	
		(8) AROR (9)	
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		bCB,2	
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Vam	_	ions in Ballons in State of St	
Sampler Name:	Preserved	Sample Intact: Y Sample ID RC-ER- 48-0799 RC-ER- 34-0799 RC-EC- 12-0799 RC-EC- 71-0799 RC-EC- 71-0799 RC-EC- 35-0799 RC-EC- 00-0799A Rinsate Blank Trip Blank Instructions: Plea	
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Submitted by:_ Agent of: Submitted by:_ Agent of: Submitted by: BDP Agent of:

Received by: Date / Time:

Agent of:

Agent of: Reliable Date / Time: 1/21/45 0400

Received by:

Agent of: Date / Time: Received by:

Reduced Standard Deliverables:

Report to: _

AHA File name: Chain of Custody for Raccoon Creek